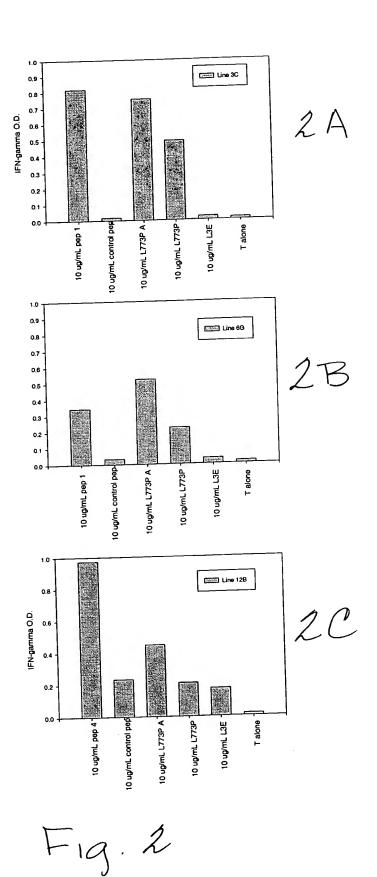
L773P Peptides	
MWQPLFFKWLLSCCPGSSQI	1-20
FFKWLLSCCPGSSQIAAAAS	6-25
LSCCPGSSQIAAAASTQPED	11-30
GSSQIAAAASTQPEDDINTQ	16-35
AAAASTQPEDDINTQRKKSQ	21-40
TOPEDDINTORKKSOEKMRE	26-45
DINTQRKKSQEKMREVTDSP	31-50
RKKSQEKMREVTDSPGRPRE	36-55
EKMREVTDSPGRPRELTIPQ	41-60
VTDSPGRPRELTIPQTSSHG	46-65
GRPRELTIPQTSSHGANRF	51-69

Fig. 1

L



D45 L773 CD4 Assay IFN-gamma Si									
	\$1 0.7 0.9 0.8 1.2 0.8 0.8 1.2 1.4	SI 0.1 1.6 0.9 1.2 0.6 0.9 1.2 1.2	Peptide SI 0.7 2.0 6.4 0.8 0.8 1.3 0.9 0.6	1 SI 1.2 1.4 0.7 0.5 0.8 1.1 0.5 0.8	Si 0.5 0.8 1.1 1.9 1.1 1.3 1.0 0.6	SI 1.2 1.3 0.7 0.8 2.0 0.7 4.2 2.4			
	0.9 2.4 0.7 0.8 0.6 1.0 0.9 2.1	1.1 1.0 0.9 1.0 2.1 1.3 1.0	Peptide 0.8 1.4 0.9 1.1 0.9 1.1 1.0	0.5 0.8 1.4 1.2 1.5 1.6 1.0	2.7 0.9 1.0 0.9 0.9 1.0 1.1	1.1 1.3 1.6 8.4 1.5 1.2 0.8 1.5			
	1.0 1.1 1.3 1.3 1.5 0.8 0.7 1.0	0.9 1.1 1.1 0.4 0.6 1.5 1.1 2.5	Peptid 1.5 1.0 0.9 1.3 1.3 1.3 1.6 0.9	1.3 0.8 0.9 1.4 0.7 0.6 0.9 2.4	1.1 1.2 2.0 0.9 1.1 1.3 2.3	0.9 0.8 1.1 1.1 0.9 1.0 0.5 0.9			
	0.9 0.9 1.0 0.9 0.8 1.1 1.1	13.6 3.9 0.9 0.8 2.6 0.8 1.0	Peptic 0.6 1.2 0.9 0.7 0.9 1.0 1.1	0.8 0.9 0.7 1.1 1.0 1.2 1.0	0.9 1.5 0.5 0.9 1.2 0.7 1.0	1:0 13.7 1.1 12.3 4.3 1.2 1.2 1.1			
	1.2 1.0 0.9 0.9 0.7 0.6 1.0	0.7 0.9 0.9 1.7 0.8 1.0 0.8	Pepti 1.0 0.9 0.8 1.1 0.9 1.1 1.1	0.5 0.7 1.2 1.4 1.3 0.5 1.0 1.0	2.1 0.7 1.3 1.5 1.4 0.8 1.2 0.8	1.2 0.8 1.5 1.5 1.3 3.1 0.6 1.7			
	0.8 1.1 1.1 0.8 0.7 1.3 1.1	0.8 0.8 0.8 1.1 1.5 1.0		0.8 0.8 0.8 0.1.1 1.1 1.0	1.0 1.0 0.7 1.1 1.1 1.4 1.6 0.8	1.5 1.1 2.5 1.2 0.7 1.1 1.8 1.2			
	0.9 0.8 1.2 1.3 2.2 1.3 1.0	1.3 2.2 1.2 1.4 0.9 1.0 1.7 0.8		0.6 0.8 1.4 0.9 1.7 1.6 1.0	1.0 1.6 0.8 1.3 0.6 1.0 1.5	1.3 0.9 1.1 0.8 0.7 1.2 1.2			
	0.8 0.9 0.7 1.6 0.9 1.1 2.5 0.6	1.6 0.8 1.0 1.0 1.0 1.0 1.1	Per 1.0 1.1 0.8 1.1 1.4 0.9 1.1	0.7 0.9 1.2 1.3 1.0 0.8 0.6 1.0	1.0 1.1 1.6 1.1 0.9 1.1 1.4	1.0 1.0 1.2 0.9 0.9 1.0 1.1			
	0.9 1.0 0.9 1.0 0.6 1.1 0.7	0.9 1.1 0.7 1.0 0.9 0.8	P4 4.9 24.4 26.3 28.1 1.3 40.2 44.0 5.1	0.7 1.2 1.2 1.1 0.7 0.9	0.8 1.3 1.8 0.8 1.2 1.3 1.1	1.3 1.3 1.2 1.0 1.1 1.2 0.9			
	0.9 1.0 0.4 0.8 1.0 1.1	1.0 1 0.9 9 0.7 0 1.0 0 1.2 1 1.0 2 0.8	1.2 1.2 0.8 0.6 1.1 1.1	1.1 0.9 1.0 1.0 1.1 1.3 0.6 1.3	0.9 0.5 1.0 1.1 0.8 1.3	1.3 1.0	3		
	1 1	1 1.2 9 0.5 1 0.5	P 2 1.2 9 1.0 7 1.4 9 1.5 9 1.7 1 1.1 8 1.3	eptide 11 1.2 1.2 1.2 1.5 1.0 7 0.9 1 1.5 3 1.0	1.6 1.1 1.4 1.0 0.5 0.5 1.6	1.5 1.5 1.9 1.9 1.0 1.1 1.1	3 5 4 1 8		

Fig. 3